

K-Nearest Neighbors

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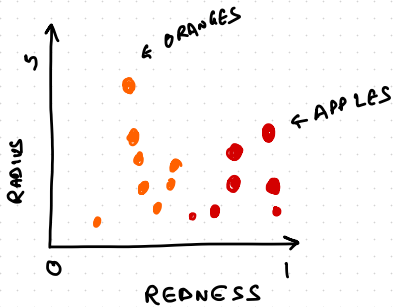
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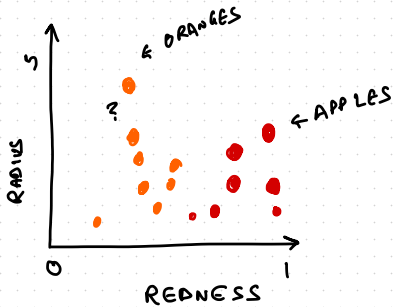
CLASSIFICATION



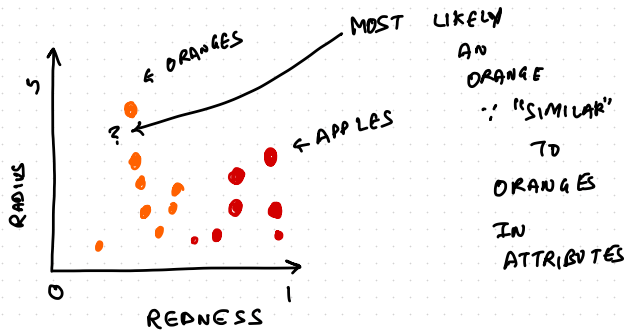
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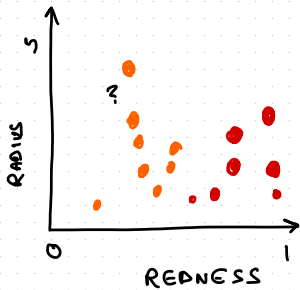
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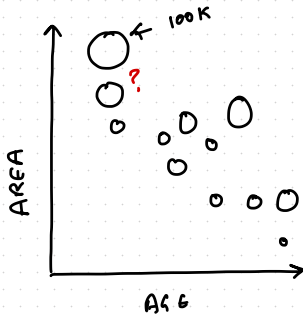
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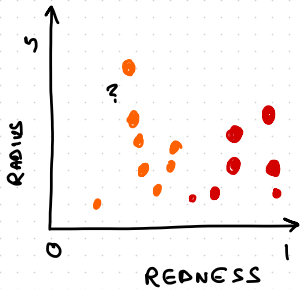
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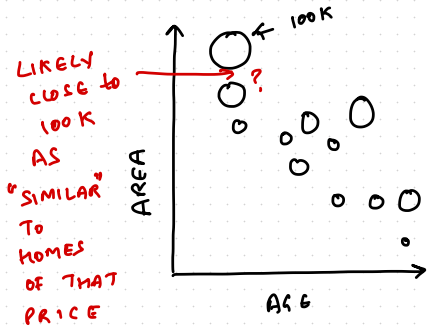
REGRESSION



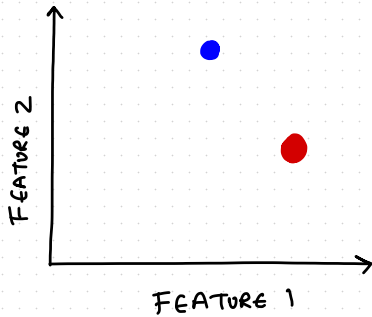
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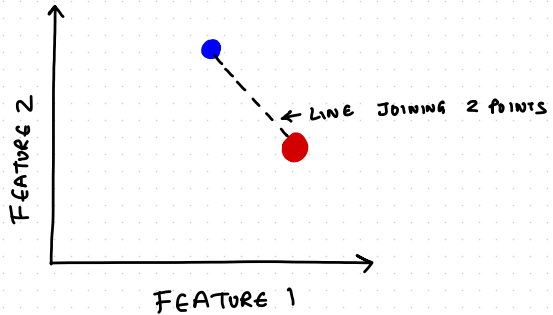
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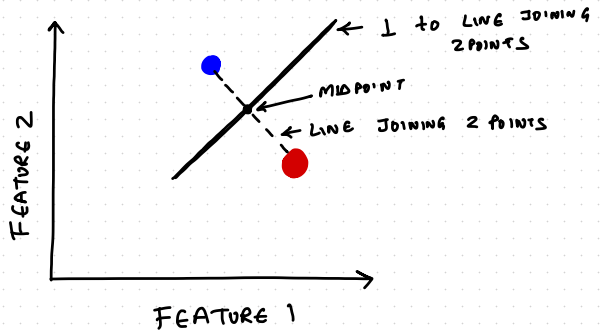
VORONOI DIAGRAM FOR 1-NN



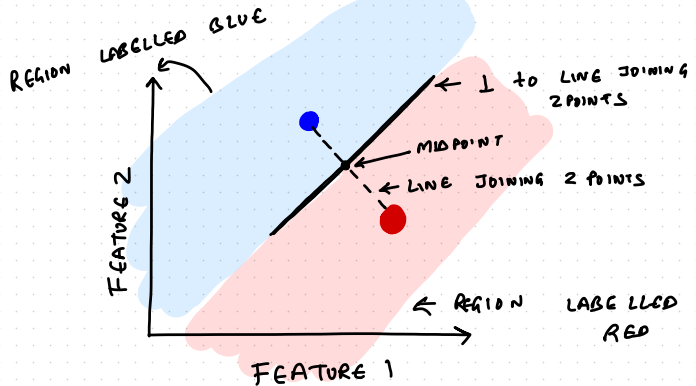
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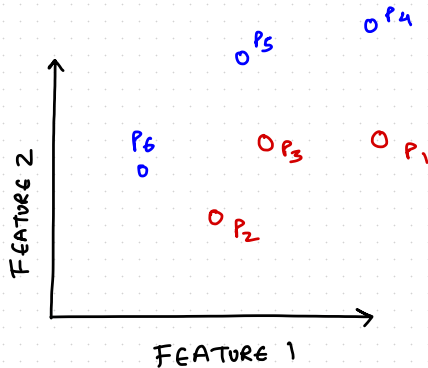
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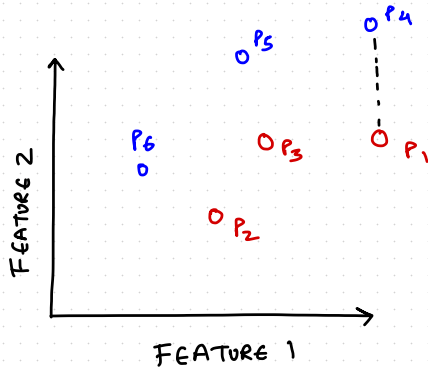
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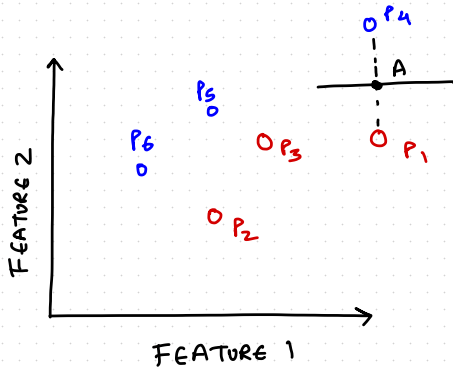


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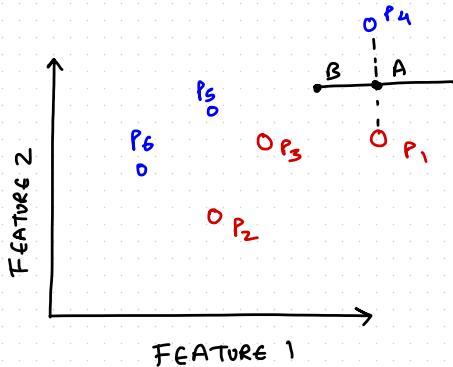
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A: MID PT B/W P_1 & P_4



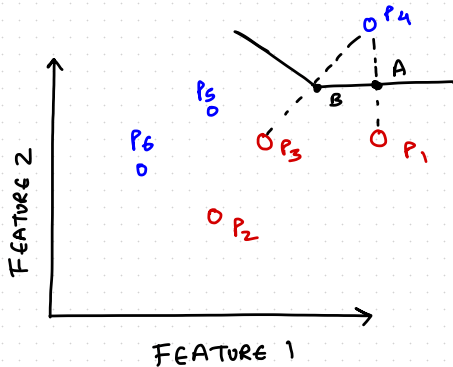
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A: MID PT B/W P_1 & P_4
B: CLOSER TO P_3 than P_1



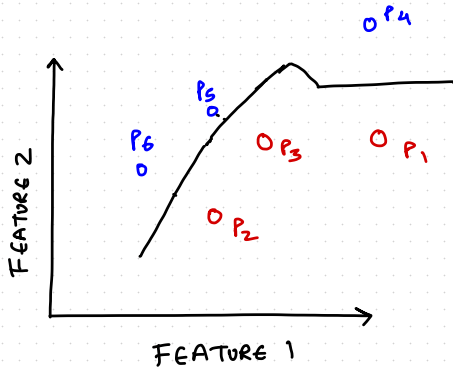
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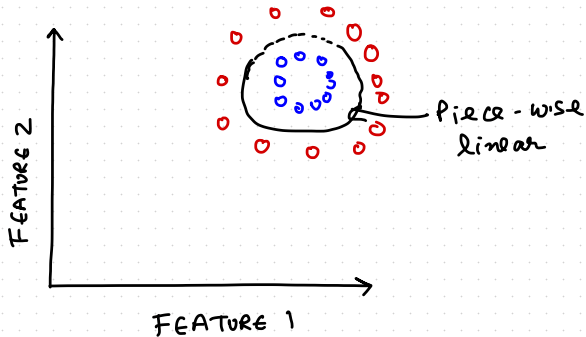
VORONOI DIAGRAM FOR 1-NN

DECISION
BOUNDARY IS
PIECE-WISE
LINEAR

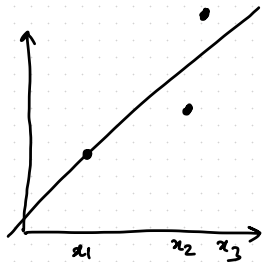


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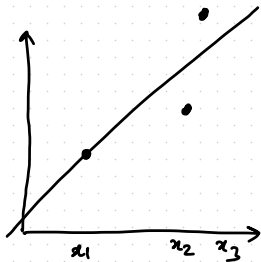


LINEAR REGRESSION

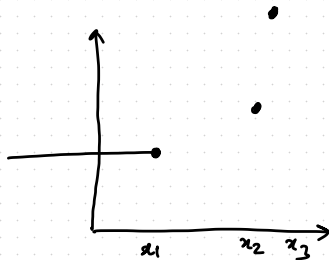


1NN REGRESSION

LINEAR REGRESSION

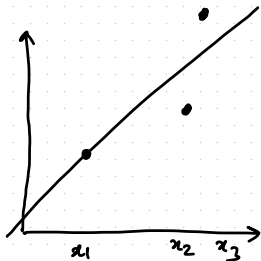


1NN REGRESSION

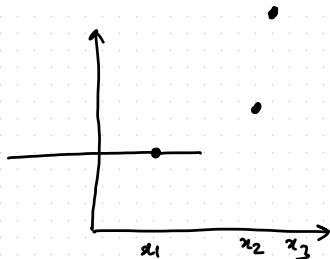


$x < x_1$: NN is (x_1, y_1)

LINEAR REGRESSION



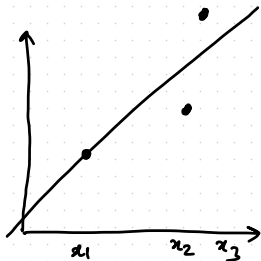
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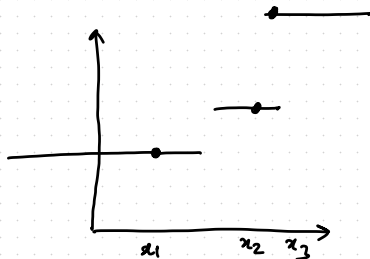
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$x < \frac{x_1 + x_2}{2}$: NN is (x_1, y_1)

LINEAR REGRESSION



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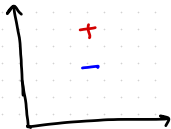


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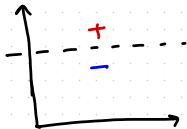
$\frac{x_1 + x_2}{2} < x < \frac{x_2 + x_3}{2}$: NN is (x_2, y_2)

KNN IS NON-PARAMETRIC



LINEAR MODEL

KNN IS NON-PARAMETRIC

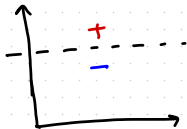


DECS
BOUNDARY

LINEAR MODEL

$$y = mx + c \quad (\# \text{ params} = 2)$$

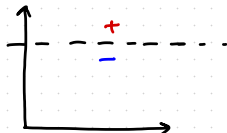
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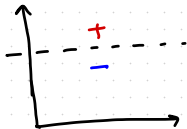
DECISION
BOUNDARY



KNN (K=1)

DECISION
BOUNDARY (LIKE $y = mx + c$)

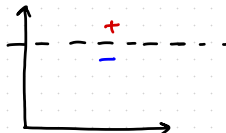
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DECISION
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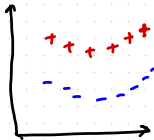


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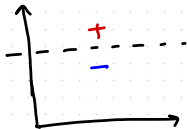
DECISION
BOUNDARY

(LIKE $y = mx + c$)

ADD DATA

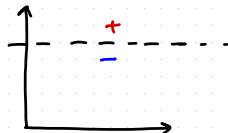


KNN IS NON-PARAMETRIC



DECISION
BOUNDARY

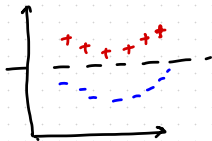
LINEAR MODEL
 $y = mx + c$ (# params = 2)



KNN (K=1)

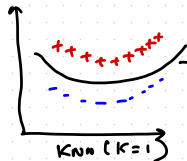
DECISION
BOUNDARY (LIKE $y = mx + c$)

ADD DATA



DECISION
BOUNDARY

LINEAR MODEL
 $y = mx + c$ (2 params)



KNN (K=1)

PARAMS $\gg 2$ (AT LEAST 10,311)

PARAMETRIC

#PARAMS FIXED
WRT DATASET SIZE

MAKE ASSUMPTIONS
(LIKE FUNCTIONAL FORM)

USUALLY QUICKER

Eg: LINEAR MODELS,
SVM (LINEAR, POLYNOMIAL)

NON-PARAMETRIC

#PARAMS GROWS
WRT DATASET SIZE

LESSER ASSUMPTIONS

USUALLY SLOWER

Eg: KNN, DT,
SVM (with
RBF)

Parametric vs Non-Parametric Models

| | Parametric | Non-Parametric |
|-------------|--|--|
| Parameter | Number of parameters is fixed w.r.t dataset size | Number of parameters grows w.r.t. to an increase in dataset size |
| Speed | Quicker (as the number of parameters are less) | Longer (as number of parameters are less) |
| Assumptions | Strong Assumptions (like linearity in Linear Regression) | Very few (sometimes no) assumptions |
| Examples | Linear Regression | KNN, Decision Tree |

Lazy vs Eager Strategies

| | Lazy | Eager |
|------------|--|---|
| Train Time | 0 | $\neq 0$ |
| Test | Long (due to comparison with train data) | Quick (as only "parameters" are involved) |
| Memory | Store/Memorise entire data | Store only learnt parameters |
| Utility | Useful for online settings | |
| Examples | KNN | Linear Regression, Decision Tree |

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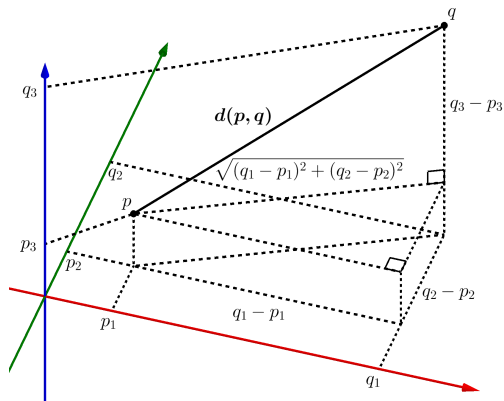
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- What is the **aggregation function** that is going to be used?
- What are the **number of neighbors** that you are going to take into consideration?
- What is the **computational complexity** of the algorithm that you are implementing?

Important Considerations: Distance Metric

The Distance Metric acts as a *measure of similarity* between the points.

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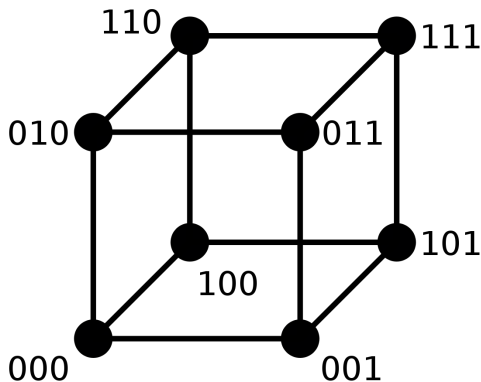
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Euclidean Distance

Important Considerations: Distance Metric

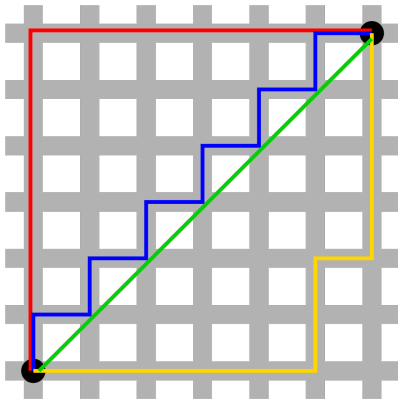
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Hamming Distance

Important Considerations: Distance Metric

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Manhattan Distance

Important Considerations: Value of K

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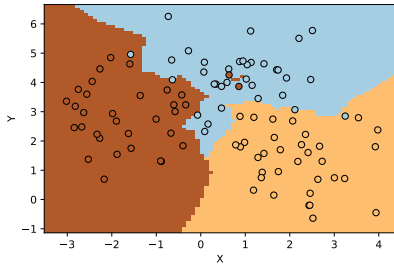
Low values of K will result in each point having a very high influence on the final output \implies noise will influence the result

High values of K will result in smoother decision boundaries

\implies lower variance but also higher bias

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$K = 3$

Important Considerations: Value of K

Aggregating data

There are different ways to go about aggregating the data from the K nearest neighbors.

- Median

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- Mode

KNN Algorithm

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 2. Predict y^*

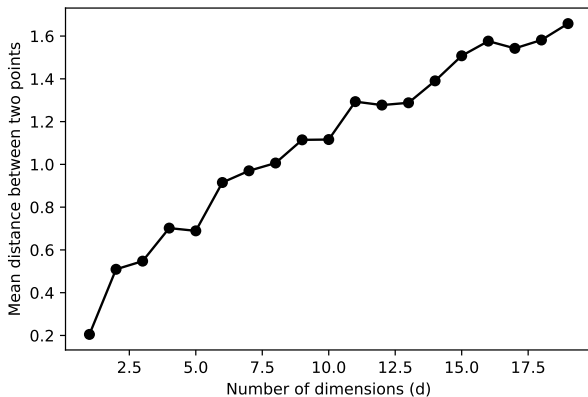
Curse of Dimensionality

With an increase in the number of dimensions:

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1. the distance between points starts to increase



For a uniformly random dataset

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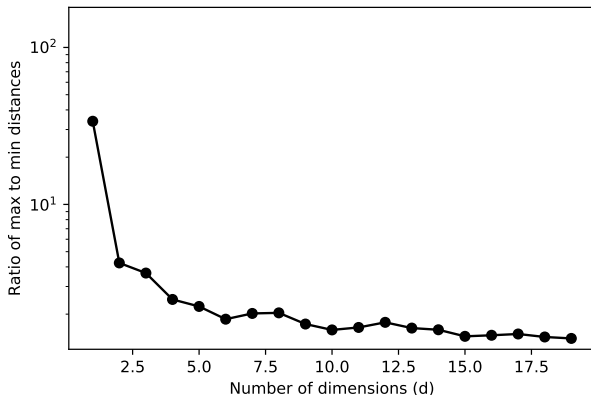
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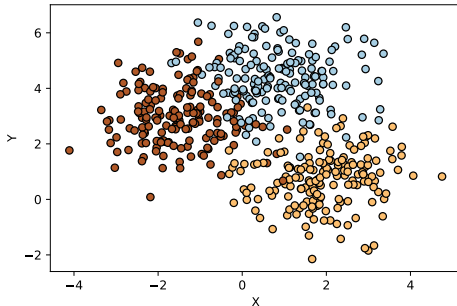
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Approximate Nearest Neighbors

Doing an exhaustive search over all the points is time consuming, especially if you have a large number of data points.



Example of a big dataset

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Such techniques include:

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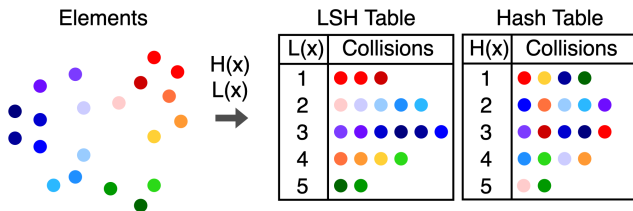
If you are willing to sacrifice accuracy there are algorithms that can give you improvements that go into orders of magnitude.

Such techniques include:

- Locality sensitive hashing
- Vector approximation files
- Greedy search in proximity neighborhood graphs

Locality sensitive hashing

Normal hash functions $H(x)$ try to keep the collision of points across bins uniform.

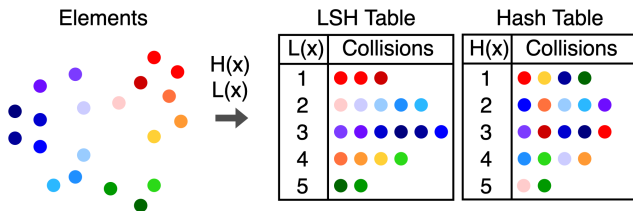


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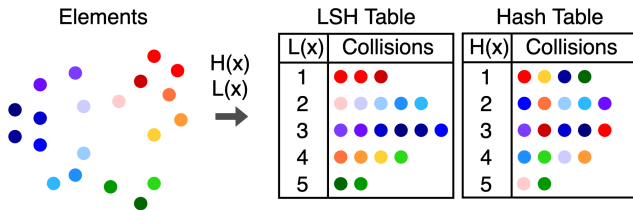


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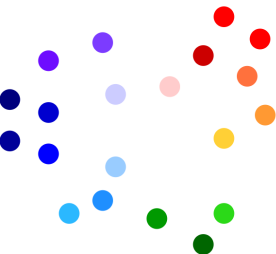
A locality sensitive hash (LSH) function $L(x)$ would be designed such that similar values are mapped to similar bins.

For such cases, all elements in a bin would be given the same label, which again can be decided on the basis of different aggregation methods



Example of a big dataset

Elements



$H(x)$
 $L(x)$



LSH Table

| L(x) | Collisions |
|------|------------|
| 1 | |
| 2 | |
| 3 | |
| 4 | |
| 5 | |

Hash Table

| H(x) | Collisions |
|------|------------|
| 1 | |
| 2 | |
| 3 | |
| 4 | |
| 5 | |

Pop Quiz: KNN Concepts

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2. Why is feature scaling important for KNN?
3. In which scenarios would you prefer KNN over parametric methods?
4. What is the time complexity of finding k nearest neighbors naively?

Key Takeaways

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- **Distance Metrics:** Choice affects performance significantly
- **Curse of Dimensionality:** Performance degrades in high dimensions
- **Scalability:** Approximate methods needed for large datasets